

European and Chinese Cooperation on Grid





Towards "Chemical" Grids

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1. Chemical Programming

2. Chemical Grid Programming

3. Perspectives







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- Initial work from Jean-Pierre Banâtre and Daniel Le Métayer (1986)
- Programming model using chemistry as a metaphor:
 - data = molecules
 - computation = chemical reactions
- Last development: HOCL





- HOCL: Higher Order Chemical Language
- \bullet Based on the $\gamma\text{-calculus}$
- A HOCL program is a chemical solution of atoms $\langle A_1, \ldots, A_n \rangle$
- Atoms A_i may be:
 - Integers, strings, ... any external object
 - Tuples A_1 :···· A_k
 - Sub-solutions
 - One-shot rules: one P by M if C
 - N-shot rules: replace P by M if C









Chemical programming : HOCL

























































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 - Conventional approaches seem to fail
 - Very high level parallel language: no process, provides atomicity and mutual exclusion, shared memory, etc.
 - No central control (locality property)
 - Chemical metaphor
 - Autonomic programs...







$$\langle \text{sieve}, 2, 3, 4, 5, 6, 7, 8, 9, 10 \rangle$$

 $\downarrow *$
 $\langle \text{sieve}, 2, 3, 5, 7 \rangle$

Stabilizing

Second Echogrid Workshop







$$\begin{array}{l} \langle \text{sieve}, 2, 3, 4, 5, 6, 7, 8, 9, 10 \rangle \\ \downarrow \ast \\ \langle \text{sieve}, 2, 3, 5, 7 \rangle \end{array}$$

Perturbation

 $\langle \text{sieve}, 2, 3, 5, 7, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20 \rangle$







$$\langle \text{sieve}, 2, 3, 4, 5, 6, 7, 8, 9, 10 \rangle$$

 $\downarrow *$
 $\langle \text{sieve}, 2, 3, 5, 7 \rangle$

Perturbation

 $\langle \text{sieve}, 2, 3, 5, 7, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20 \rangle$ *Re-stabilizing* $\downarrow *$ $\langle \text{sieve}, 2, 3, 5, 7, 11, 13, 17, 19 \rangle$





Application level: write applications with HOCL

System level: specification and interface of "chemical" grids with HOCL



 Applications are chemical programs

Application level

- No reference to any grid mechanism
- HOCL used as a coordination language







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import Pixel, Ray, LightRay, Contrib, Scene.





import Pixel, Ray, LightRay, Contrib, Scene. $\langle p_1, \ldots, p_n, \text{ renderPixel} \rangle$





 ${\bf import}\, {\rm Pixel}, {\rm Ray}, {\rm LightRay}, {\rm Contrib}, {\rm Scene}.$

 $\langle p_1, \ldots, p_n, \text{ renderPixel} \rangle$

 $\begin{aligned} \text{renderPixel} = \textbf{replace } p :: \text{Pixel} \\ \textbf{by } & \langle firstRay(p), \text{ renderRay, deleteRay,} \\ & \text{enlighten, sumContrib} \end{aligned}$





import Pixel, Ray, LightRay, Contrib, Scene. $\langle p_1, \ldots, p_n, \text{ renderPixel} \rangle$ renderPixel = replace p::Pixel **by** $\langle firstRay(p), renderRay, deleteRay,$ enlighten, sumContrib> renderRay = replace r::Ray by Scene. intersectRays(r)if $r.contribution() > \epsilon$ deleteRay = replace r::Ray, ω by ω if r.contribution() $\leq \epsilon$ enlighten = replace l::LightRay **by** *l.computeContrib()* sumContrib = replace c_1 ::Contrib, c_2 ::Contrib **by** $c_1.add(c_2)$



System level



- Grid viewed as a chemical solution:
 - resources = molecules
 - coordination = chemical reactions
- Chemical program as:
 - a specification
 - an interface for the grid administrator



Credit: Andrew Wood, Nick Drew, Russell Beale and

Bob Hendley, 1995.





 $\langle R_1:\langle \langle r_1 \rangle, \ldots, \langle r_n, \rangle \rangle, R_2:\langle \rangle, \ldots, R_n:\langle \rangle, renderRules, split, merge, newRes, remRes \rangle$





 $\langle R_1: \langle \langle r_1 \rangle, \dots, \langle r_n, \rangle \rangle, R_2: \langle \rangle, \dots, R_n: \langle \rangle, renderRules, split, merge, newRes, remRes \rangle$

split = replace $res_1:\langle\langle r::Ray, \omega_p \rangle, \omega_1 \rangle, res_2:\langle\omega_2 \rangle$ by $res_1:\langle\omega_1 \rangle, res_2:\langle\langle r, \omega_p \rangle, \omega_2 \rangle$ if $res_1.load() >> res_2.load()$





 $\langle R_1: \langle \langle r_1 \rangle, \dots, \langle r_n, \rangle \rangle, R_2: \langle \rangle, \dots, R_n: \langle \rangle, renderRules, split, merge, newRes, remRes \rangle$

split = **replace**
$$res_1:\langle\langle r::Ray, \omega_p \rangle, \omega_1 \rangle, res_2:\langle\omega_2 \rangle$$

by $res_1:\langle\omega_1 \rangle, res_2:\langle\langle r, \omega_p \rangle, \omega_2 \rangle$
if $res_1.load() >> res_2.load()$

merge = **replace**
$$res_1:\langle \omega_1 \rangle$$
, $res_2:\langle \langle c::Contrib \rangle$, $\omega_2 \rangle$
by $res_1:\langle \omega_1, \langle c \rangle \rangle$, $res_2:\langle \omega_2 \rangle$
if $res_1 < res_2$







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- Currently EchoGrid Fellow at the ICT
- Initial work:

Application level: implement HOCL using GSML





谢谢

Thank You!

Second Echogrid Workshop

